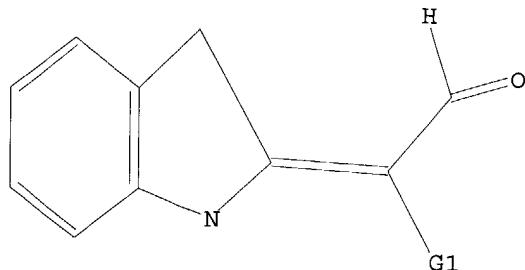


Uploading C:\Program Files\Stnexp\Queries\10055664.str

L1 STRUCTURE UPLOADED

=> d  
L1 HAS NO ANSWERS  
L1 STR



G1 CH<sub>2</sub>O

MP search

Structure attributes must be viewed using STN Express query preparation.

=> s 11  
SAMPLE SEARCH INITIATED 10:15:33 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 59 TO ITERATE  
  
100.0% PROCESSED 59 ITERATIONS 0 ANSWERS  
SEARCH TIME: 00.00.01  
  
FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 720 TO 1640  
PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s 11 full  
FULL SEARCH INITIATED 10:15:37 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 1283 TO ITERATE

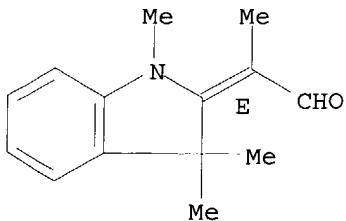
100.0% PROCESSED 1283 ITERATIONS 2 ANSWERS  
SEARCH TIME: 00.00.01

L3 2 SEA SSS FUL L1

=> d 1-2

L3 ANSWER 1 OF 2 REGISTRY COPYRIGHT 2004 ACS on STN  
RN 124244-70-8 REGISTRY  
CN Propanal, 2-(1,3-dihydro-1,3,3-trimethyl-2H-indol-2-ylidene)-, (E)- (9CI)  
(CA INDEX NAME)  
FS STEREOSEARCH  
MF C14 H17 N O  
SR CA  
LC STN Files: BEILSTEIN\*, CA, CAPLUS  
(\*File contains numerically searchable property data)  
DT.CA CAPplus document type: Journal  
RL.NP Roles from non-patents: PREP (Preparation); RACT (Reactant or reagent)

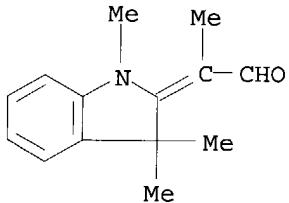
Double bond geometry as shown.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

1 REFERENCES IN FILE CA (1907 TO DATE)  
 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L3 ANSWER 2 OF 2 REGISTRY COPYRIGHT 2004 ACS on STN  
 RN 42909-52-4 REGISTRY  
 CN Propanal, 2-(1,3-dihydro-1,3,3-trimethyl-2H-indol-2-ylidene)- (9CI) (CA INDEX NAME)  
 OTHER NAMES:  
 CN 1,3,3-Trimethyl-2-( $\alpha$ -formylethylidene)indoline  
 CN 1,3,3-Trimethyl-2-(1-formylethylidene)indoline  
 FS 3D CONCORD  
 MF C14 H17 N O  
 LC STN Files: BEILSTEIN\*, CA, CAPLUS  
     (\*File contains numerically searchable property data)  
 DT.CA CAplus document type: Patent  
 RL.P Roles from patents: PREP (Preparation); USES (Uses)



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

3 REFERENCES IN FILE CA (1907 TO DATE)  
 3 REFERENCES IN FILE CAPLUS (1907 TO DATE)

>>> FOR SEARCHING PREPARATIONS SEE HELP PRE <<<

\*\*\*\*\*  
\* PLEASE NOTE THAT THERE ARE NO FORMATS FREE OF COST. \*  
\* SET NOTICE FEATURE: THE COST ESTIMATES CALCULATED FOR SET NOTICE \*  
\* ARE BASED ON THE HIGHEST PRICE CATEGORY. THEREFORE; THESE \*  
\* ESTIMATES MAY NOT REFLECT THE ACTUAL COSTS. \*  
\* FOR PRICE INFORMATION SEE HELP COST \*  
\*\*\*\*\*

=> s 11 full  
FULL SEARCH INITIATED 10:16:14 FILE 'BEILSTEIN'  
FULL SCREEN SEARCH COMPLETED - 651 TO ITERATE

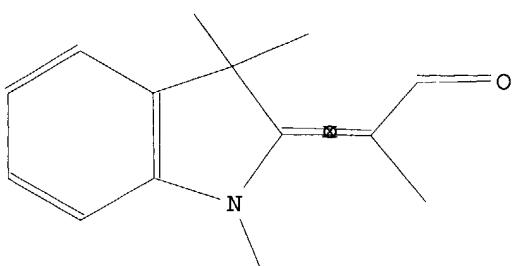
100.0% PROCESSED 651 ITERATIONS 1 ANSWERS  
SEARCH TIME: 00.00.10

L4 1 SEA SSS FUL L1

=> d ide

L4 ANSWER 1 OF 1 BEILSTEIN COPYRIGHT 2004 BEILSTEIN MDL on STN

Beilstein Records (BRN) : 3545115  
Beilstein Pref. RN (BPR) : 124244-70-8  
CAS Reg. No. (RN) : 42909-52-4, 124244-70-8  
Chemical Name (CN) : 2-(1,3,3-Trimethyl-2-indolinyliden)propanal  
Autonom Name (AUN) : 2-(1,3,3-trimethyl-1,3-dihydro-indol-2-ylidene)-propionaldehyde  
Molec. Formula (MF) : C14 H17 N O  
Molecular Weight (MW) : 215.29  
Lawson Number (LN) : 25479, 2817  
File Segment (FS) : Stereo compound  
Compound Type (CTYPE) : heterocyclic  
Constitution ID (CONSID) : 3137287  
Tautomer ID (TAUTID) : 3361259  
Beilstein Citation (BSO) : 6-21  
Entry Date (DED) : 1991/10/23  
Update Date (DUPD) : 1991/11/01



Field Availability:

Code	Name	Occurrence
BRN	Beilstein Records	1

BPR	Beilstein Preferred RN	1
RN	CAS Registry Number	2
CN	Chemical Name	1
AUN	Autonomname	1
MF	Molecular Formula	1
FW	Formular Weight	1
LN	Lawson Number	2
FS	File Segment	1
CTYPE	Compound Type	1
CONSID	Constitution ID	1
TAUTID	Tautomer ID	1
BSO	Beilstein Citation	1
ED	Entry Date	1
UPD	Update Date	1
MP	Melting Point	1
NMR	Nuclear Magnetic Resonance	3
UVS	UV and Visible Spectrum	1

This substance also occurs in Reaction Documents:

Code	Name	Occurrence
RX	Reaction Documents	3
RXREA	Substance is Reaction Reactant	2
RXPRO	Substance is Reaction Product	1

=>

Connection closed by remote host